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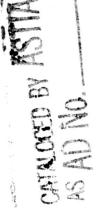
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MEMORANDUM RM-3084-PR APRIL 1962



POLYNOMIAL APPROXIMATION—A NEW COMPUTATIONAL TECHNIQUE IN DYNAMIC PROGRAMMING—I: ALLOCATION PROCESSES

Richard Bellman, Robert Kalaba and Bella Kotkin

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PREFACE

Part of the Project RAND research program consists of basic supporting studies in mathematics. One aspect of this is concerned with optimization processes. In this field, the technique of dynamic programming has developed into a powerful mathematical tool.

In the present Memorandum the authors investigate the applicability of polynomial approximation as an adjunct to the technique of dynamic programming.

SUMMARY

In this Memorandum the authors initiate the application of the simple yet powerful computational technique of polynomial approximation to problems in dynamic programming.

The theoretical applicability of orthogonal polynomials is first discussed and then applied to one— and two—dimensional allocation problems. Numerical results obtained from FORTRAN programs involving Legendre polynomials are presented.

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POLYNOMIAL APPROXIMATION—A NEW COMPUTATIONAL TECHNIQUE IN DYNAMIC PROGRAMMING—I: ALLOCATION PROCESSES

1. INTRODUCTION

The problem of maximizing the function

(1.1)
$$F(x_1, x_2, ..., x_N) = g_1(x_1) + g_2(x_2) + \cdots + g_N(x_N)$$

over the domain

(1.2)
$$x_1 + x_2 + \cdots + x_N = x, \quad x_1 \ge 0$$

can be reduced via familiar dynamic programming techniques (see [1]) to that of determining the sequence of functions $\{f_n(x)\}$ generated by means of the recurrence relation

(1.3)
$$f_{1}(x) = g_{1}(x),$$

$$f_{n+1}(x) = \max_{0 \le y \le x} [g_{n+1}(y) + f_{n}(x - y)].$$

The problem can thus be solved numerically in a very simple fashion, regardless of the complexity of the functions $g_1(x)$. A number of important allocation processes can be resolved in this way. If we consider cases in which two distinct types of resources must be allocated, we face the problem of maximizing the function

(1.4)
$$F(x_1, x_2, ..., x_N; y_1, y_2, ..., y_N)$$

$$= g_1(x_1, y_1) + g_2(x_2, y_2) + ... + g_N(x_N, y_N)$$

over the domain

(1.5)
$$x_1 + x_2 + \cdots + x_N = x, \quad x_1 \ge 0,$$

 $y_1 + y_2 + \cdots + y_N = y, \quad y_1 \ge 0.$

Theoretically, there is no difficulty in applying the same techniques.

The maximization problem can be reduced to the determination of the sequence $\{f_n(x,y)\}$ generated by means of the recurrence relation

(1.6)
$$f_{n+1}(x,y) = \max_{\substack{0 \le w \le x \\ 0 \le r \le y}} [g_{n+1}(w,r) + f_n(x - w,y - r)].$$

In principle, this equation can be solved computationally using the same technique that applies so well to (1.3). In practice (see [1] for a discussion), questions of time and accuracy arise. There are a number of ways of circumventing these difficulties, among which the Lagrange multiplier plays a significant role.

In this series of papers, we wish to present a number of applications of a new, simple and quite powerful method, that of polynomial approximation. We shall begin with a discussion of the allocation process posed in the foregoing paragraphs and continue, in subsequent papers, with a treatment of realistic trajectory and guidance processes. In a separate series of papers we shall apply this fundamental attack upon

dimensionality to the solution of a number of the equations of mathematical physics.

We would like to express our appreciation to Oliver Gross for the analytic solution of some test problems we used to check the accuracy of our techniques, and for his general interest in the program.

2. POLYNOMIAL APPROXIMATION

In the systematic study of dynamic programming as a computational algorithm given in [1], a function f(x) is considered to be a table of values at an appropriate set of grid points:

$$\begin{array}{c|cccc}
x & f(x) \\
0 & f_0 \\
\Delta & f_1 \\
2\Delta & f_2 \\
\vdots & \vdots \\
K\Delta & f_K
\end{array}$$

In this way, the function f(x) is stored in the computer. For functions of one variable with K=100 or 1000, this is a reasonably convenient way to proceed. For functions of two variables, this procedure becomes a bit inconvenient since (K+1) values for x combined with (K+1) values for y yields a total set of $(K+1)^2$ values. Consequently, when we encounter functions of three or more variables, we must

balance accuracy against time and the limited storage of contemporary computers in our choice of K.

The storage of functional values by means of a table of values is ideally suited to the treatment of problems involving functions of quite arbitrary form. It is, on the other hand, quite wasteful and inefficient if we are dealing with functions possessing a definite structure, which is to say, situations in which there is a high correlation between the values of $f(r\Delta)$ and $f(s\Delta)$. Since functions of this type occur in many important applications, and throughout mathematical physics, it is worthwhile to develop methods which take advantage of the "smoothness" of the function.

One such method is <u>polynomial approximation</u>, or to be precise, generalized polynomial approximation. We represent the function in the form

(2.2)
$$f(x) \stackrel{\sim}{=} \sum_{k=1}^{M} a_k \varphi_k(x),$$

where the $\varphi_k(x)$ are known elementary functions such as x^k , $\sin kx$, $P_k(x)$ (the Legendre polynomial of degree k) or $T_k(x)$ (the Cebycev polynomial of degree k), and then store the function for all values of interest by means of the set of M coefficients $[a_1, a_2, \ldots, a_M]$.

It is important to point out that (2.2) is particularly useful in automatically furnishing the interpolation—values frequently needed in dynamic programming calculations. If

one uses a table of values of the form shown in (2.1), interpolation is frequently a source of difficulty.

To determine the coefficients $a_{\bf k}$ it is convenient to make the $\phi_{\bf k}(x)$ be an orthonormal set. Then

(2.3)
$$a_k = \int_{0}^{1} f(x)\phi_k(x)dx$$

assuming for the purposes of convenience that $0 \le x \le 1$. We can, of course, use a Cebycev fit instead, and we will explore this in subsequent papers. A priori, we would suspect that it would be more efficient to use a mean-square fit (implied by (2.3)), and take M to be larger, than to go to the trouble of determining the a_k to minimize the function

(2.4)
$$\max_{0 \le x \le 1} |f(x) - \sum_{k=1}^{M} a_k \varphi_k(x)|,$$

which for a fixed M may be expected to yield a more accurate approximation. Alternatively, one could use an approximation by polygonal functions [1].

To evaluate the a_k without requiring a knowledge of too many values of f(x), we use a quadrature technique

(2.5)
$$a_{k} \stackrel{\cong}{=} \sum_{1=1}^{R} w_{1}f(x_{1})\varphi_{k}(\overline{x}_{1}),$$

where the weights $\mathbf{w_i}$ and the quadrature points $\overline{\mathbf{x_i}}$ are chosen so that the equation

(2.6)
$$\int_{0}^{1} g(x)dx = \sum_{i=1}^{R} w_{i}g(\overline{x}_{i})$$

is exact for polynomials in x of degree (2R-1) or less. This requirement narrows us down to the Gauss quadrature technique [2]. If we use generalized polynomials (expressions such as (2.6)), we will obtain different weights and quadrature points.

We may then store the function f(x) for $0 \le x \le 1$ by storing the coefficients a_k or the particular values $f(\overline{x}_i)$ which enable us to compute the a_k .

3. APPLICATION TO DYNAMIC PROGRAMMING

Consider now the application of these ideas to the computational solution of the functional equations of dynamic programming. Suppose that we wish to compute the sequence $\{f_n(x)\}$ determined by

(3.1)
$$f_N(x) = \max_{0 \le y \le x} [g_N(y) + f_{N-1}(x - y)],$$

 $N \geq 2$, given that $f_1(x) = g_1(x)$. To avoid the tabulation of each of the functions $f_N(x)$ at the x-grid $[0,\Delta,\ldots,K\Delta]$, where K may be a large number, we approximate to each function $f_N(x)$ in the manner indicated above. Starting with $f_1(x)$, we store the values $f_1(x_1)$, $i=1,2,\ldots,R$, needed to evaluate $f_1(x-y)$ in the formula determining $f_2(x)$,

(3.2)
$$f_2(x) = \max_{0 \le y \le x} [g_2(x_2) + f_1(x - y)].$$

We then determine successively $f_2(x_1)$, $f_2(x_2)$,..., $f_2(x_R)$, a set of values which stores the function $f_2(x)$. This process is then repeated.

Although the calculation of $f_{n-1}(x-y)$ using (2.2) is far more time-consuming than taking a value from storage, we expect to gain time over-all because of the fact that we are required to calculate only a few values, $f_n(x_i)$, $i=1,2,\ldots,R$, at each stage.

4. THE LEGENDRE POLYNOMIALS

Since in allocation processes we have a range $[0,x_0]$ which stays constant as the process continues from stage to stage, we can normalize and consider the basic interval to be [0,1]. Since the interval is finite and we want, at the moment, a polynomial approximation, we shall employ Legendre polynomials.

(4.1)
$$\varphi_k(x) = (2k + 1)^{1/2} P_k(2x - 1).$$

The sequence $\{\phi_k(x)\}$ is then orthonormal over [0,1], i.e.,

(4.2)
$$\int_{O}^{1} \varphi_{k}(x) \varphi_{\ell}(x) dx = \delta_{k\ell}.$$

From the standard recurrence relations for $P_{\mathbf{k}}(\mathbf{x})$, we obtain the relation

$$\phi_{1}(x) = 1, \quad \phi_{2}(x) = 3^{1/2}(2x - 1),$$

$$\phi_{1+2}(x) = \frac{(2i + 3)^{1/2}}{(i + 1)} \left[(2i + 1)^{1/2}(2x - 1)\phi_{i+1}(x) - \frac{i}{(2i - 1)^{1/2}} \phi_{i}(x) \right].$$

This relation makes the evaluation of the sequence of values of $\phi_{\bf i}(x)$ for any x a relatively simple matter.

o. EXAMPLES

In order to experiment with this new approximating procedure, we devised a FORTRAN program for the IBM-7090 whereby at each stage, after obtaining the new coefficients $a_k^{(1)}$, we computed $f_i(x)$ from (2.2) for a succession of as many values of x as desired, and printed the result after each computation. We used two modes of output, either a list of numerical values, $[x,f_i(x)]$, or a graphical plot (done directly by the 7090) of x versus $f_i(x)$.

We experimented with several types of functions $g_1(x)$ for which the results could be derived from analytic considerations. Using the known analytic results as a checkpoint, we varied the parameters R, M, and the grid size H, in order to determine the degree of accuracy we might expect in general. We found remarkably good agreement to 2

 $^{\ \ ^{\}circ}$ Oliver Gross was of considerable assistance to us in this respect.

or more significant figures with a relatively low order of approximation, namely R = 5, M = 6. This is encouraging from the point of view of extending the method in the experimental investigations of higher—dimensional allocation processes where, as pointed out above, time and storage aspects become significant. We also incorporated in our program restraints of the form $0 \le a_1 \le x_1 \le b_1 \le x$, since constraints of this nature often occur in applications.

a. Time Estimates

Following are some estimates of the execution time required on the 7090:

- 1. 10 seconds for 4 stages, R = 5, M = 6 and a grid of 0.05 both for the search and output listing.
- 2. 3 minutes for a total of 4 cases of 10 stages each; R = 3, M = 4; R = 5, M = 6; R = 7, M = 8; R = 10, M = 11. The grid in the search for the maximum is 0.01 and the output listing is given for every x along the grid.
- 3. 3 minutes for the total of 4 cases mentioned above, 10 stages per case and a grid of 0.01 for the search. The output is a graph where the independent variable x is listed at intervals of 0.025.

b. Numerical Results

1. $g_1(x) = i(x)^{1/2}$. Using the Schwarz inequality, we readily obtain the values

$$f_N(x) = (\frac{N}{6}(N+1)(2N+1)x)^{1/2}.$$

For R = 10, M = 11, H = 0.01 we found poor agreement at the origin. This is to be expected because of the infinite slope at x = 0. Agreement between exact and computed values was good as soon as x moved away from the singular value 0, as can be seen from the following table:

Function	Exact	Computed
f ₁ (0)	0.0	0.064
f ₁ (2)	0.448	0.447
f ₁ (1)	1.00	1.00
f ₁₀ (0)	0.0	3.13
f ₁₀ (1)	19.6	19.6

- 2. $g_1(x) = i(x + 0.1)^{1/2}$. We avoided the previous difficulty at x = 0 (see the computed value of $f_{10}(0)$ above), but found the function still rather sensitive near the origin. As N (the number of stages) increased, the agreement at x = 0 decreased.
- 3. $g_1(x) = 1(x + 1)^{1/2}$. The Schwarz inequality yields the upper bound

$$f_N(x) \le (\frac{N}{6}(N+1)(2N+1)(x+N))^{1/2}$$
.

However, since the x_1 are subject to the restraint $0 \le x_1 \le x$, we do not necessarily achieve the upper bound. Some exact values based on an analysis by 0. A. Gross are listed as check points, R = 10, M = 11, H = 0.01.

Function	Upper bound	Exact value	Computed value
f ₁ (0)	1.00	1.00	1.00
f ₁ (1)	1.41	1.41	1.41
f ₃ (0)	6.48		6.00
f ₃ (.5)	7.00	6.67	6.67
f ₇ (.35)	32.1	29.1	29.1
f ₁₀ (1)		59•3	59•3

4.
$$g_1(x) = e^{-5/(1+10x)}$$
, R = 10, M = 11, H = .01.

Function	Exact	Computed
f ₂ (.2)	0.196	0.197
f ₂ (.7)	0.659	0.659
f ₃ (.2)	0.203	0.204
r ₃ (.9)	0.860	0.862
f ₅ (.2)	0.218	0.217
f ₁₀ (1)	1.001	1.001

6. TWO-DIMENSIONAL APPROXIMATION

The problem of maximizing the function

(6.1)
$$\sum_{i=1}^{N} g_{i}(x_{i}, y_{i})$$

over the domain

(6.2)
$$\sum_{1=1}^{N} x_1 = x, \quad 0 \le a_1 \le x_1 \le b_1, \quad x_1 \le x,$$

$$\sum_{1=1}^{N} y_1 = y, \quad 0 \le c_1 \le y_1 \le d_1, \quad y_1 \le y,$$

can be readily handled by the methods described in Secs. 2 and 3.

The dynamic programming recurrence relation is:

$$f_{N}(x,y) = \max_{R} [g_{N}(x_{N},y_{N}) + f_{N-1}(x - x_{N},y - y_{N})],$$

where R is determined by $a_N \leq x_N \leq \min(x,b_N)$, $c_N \leq y_N \leq \min(y,d_N)$. Let \overline{x}_1 , $i=1,\ldots,R$, be the roots of the normalized Legendre polynomial $\emptyset_R(x)$, and let $\{\overline{y}_j\}$ be a duplicate set of these quadrature points. Each function $f_N(x,y)$ is expressed approximately by the relation

(6.3)
$$f_{N}(x,y) = \sum_{r=1}^{M} \sum_{s=1}^{M} a_{r,s}(N) \phi_{r}(x) \phi_{s}(y),$$

where the coefficients, using the quadrature method, are given by

(6.4)
$$a_{k,s}^{(N)} = \sum_{i=1}^{R} \sum_{j=1}^{R} w_i w_j f_N(\overline{x}_i, \overline{y}_j) \phi_k(\overline{x}_i, \overline{y}_j).$$

As in the one-dimensional case, we start with the known values $f_1(x,y) = g_1(x_1,y_1)$, $x_1 = x$, $y_1 = y$. In stage n we store the values $f_n(\overline{x}_1,\overline{y}_j)$, for $i,j=1,2,\ldots,M$, and then compute and store the values $a_{r,s}(n)$ in the storage allotted to the previous stage. The latter coefficients are utilized in the computation of $f_n(x-x_{n+1},y-y_{n+1})$ to obtain the values of $f_{n+1}(x_1,y_j)$ in the next stage.

7. EXAMPLES

a. Time Estimate

The execution time required on the 7090 was 2 minutes for 4 stages, with R = 5, M = 6, a grid of H = 0.05 in the two-dimensional search, and an output listing of 3 test values of f(x,y) in each stage.

b. Numerical Results

1. $g_1(x,y)=(2i-1)^{1/2}(xy)^{1/4}$. Using the Hölder inequality, we readily obtain the exact values $f_n(x,y)=n(xy)^{1/4}$, $x_n(x,y)=(2n-1)x/n^2$, $y_n(x,y)=(2n-1)y/n^2$. Our results for R=5, M=6, and a grid of 0.05 in the search are:

Function	Exact	Computed
f ₂ (.5,.5)	1.40	1.40
$x_2(.5,.5)$	0.375	0.35 (to the nearest 0.05)
f ₄ (1,1)	4.00	3.91

Here again the agreement was poor at the origin, presumably because of the singular behavior of $(xy)^{1/4}$ at x=0, y=0. To confirm this hypothesis, we considered the next case.

2. $g_i(x,y) = (x + iy)/(1 + x + iy)$. This case, as well as the theoretical values, was suggested by O. A. Gross, and he determined the exact values.

Function	Exact	Computed
f ₂ (1,1)	1.17	1.17
f ₂ (1,0)	0.667	0.647

8. DISCUSSION

As can be seen, the agreement in general is quite satis—factory. We can obtain reasonably accurate values of the maximum return and of the optimal allocation policy using small amounts of machine time.

Combining these techniques with the method of the Lagrange multiplier, we can expect to solve three-and four-dimensional resource allocation problems. Extending the method to cover the approximation of functions of 3, 4, 5 and 6 variables, we can treat Hitchcock-Koopmans allocation processes of quite high dimension.

Finally, if we combine these techniques—polynomial approximations and Lagrange multipliers—with that of successive approximations, there should be very few allocation processes which still resist our efforts.

```
Appendix A
   1 DIMENSIONAL ALLOCATION VIA DYNAMIC PROGRAMMING, APPROXIMATIONS
       LIST
       LABEL
C649501
      COPMON N, P, JR, H, IH, L, Y, XL, W, A, P, G, NI, A1, B1
     X , IH1, L1, L2, H1, S, S1, HOLAST, HOLBL, HOLDOT
      DIMENSION XL(20),W(20),A(20),P(20),F(20),X(20),A1(10),B1(10)
    1 CALL INPUT
      IF(N-100)2,3,3
C
      STAGE 1
    2 NI =1
      DO 4 K=1.P
      SUM=0.
      DC 5 J=1,JR
      Y=XL(J)
      CALL RETG
      CALL POLY
      F(J)=G
      Y=(L)X
    5 SUM=SUM+W(J)+F(J)+P(K)
      A(K)=SUP
    4 CONTINUE
      DUTPUT 42,60,NI,JR,M
   60 FURMATIZOHI
                     ITERATIVE STAGE
                                        15,5H
                                                 R=13,5H
                                                            M=13
      OUTPUT 42,661
  661 FURMAT(53HO
                       DECISION
                                            RETURN AT R ROOT VALUES
      OUTPUT 42,61,(X(J),F(J),J=1,JR)
   61 FORMAT(1H 2E20.8)
      CUTPUT 42.550
  550 FORMAT (30HO
                      COEFFICIENTS A(K)
      CUTPUT 42,50,(A(K),K=1,M)
   50 FURMAT(1H E20.8)
      CALL OUT
C
      STAGE NI GREATER THAN 1
      DO 6 N2=2.N
      NI=N2
      DO 7 J=1,JR
      Y=0.
      CALL RETG
      Y=XL(J)-0.
      CALL POLY
      SUP=0.
      DC 8 K=1.M
    8 SUM=SUM+A(K)+P(K)
      F(J)=G+SUP
      X(J)=0.
      CO 9 1=2,1H
      X1=1
      IF((XI-1.)+H-A1(N2))9,110,110
  110 CONTINUE
      IF((XI-1.) • H-61(N2))111,111,11
  111 CONTINUE
      IF((XI-1.) • H-XL(J))10,10,11
  10 Y=(XI-1.) OH
      CALL RETG
      Y=XL(J)-(XI-1.) OH
      CALL POLY
      SUM=0.
     DO 12 K=1.M
  12 SUP=SUM+A(K) +P(K)
```

)

```
PH=G+SUP
      IF(PH-F(J))9,9,15
   15 F(J)=PH
      X(J)=(XI-1.)+H
    9 CONTINUE
   11 CONTINUE
    7 CONTINUE
      CUTPUT 42,60,NI,JR,M
      CUTPUT 42.661
      OUTPUT 42,61,(X(J),F(J),J=1,JR)
      DO 16 K=1,M
      SUM=0.
      DO 17 J=1.JR
      Y=XL(J)
      CALL POLY
   17 SUM=SUM+W(J)+F(J)+P(K)
      A(K)=SUM
   16 CONTINUE
      CUTPUT 42,550
      OUTPUT 42,50,(A(K),K=1,M)
      CALL OUT
    6 CONTINUE
      GO TO 1
    3 CALL EXIT
      STOP
      END
       LIST
       LABEL
C649502
      INPUT ALL CONSTANTS
      SUBROUTINE INPUT
      COMMON N,M,JR,H,IH,L,Y,XL,W,A,P,G,NI,A1,B1
     X , IH1, L1, L2, H1, S, S1, HOLAST, HOLBL, HOLDOT
      DIMENSION XL(20),W(20),A(20),P(20),F(20),X(20),A1(10),B1(10)
      INPUT 41,20,N.M.JR.IH.L
      INPUT 41,21, (XL(J),J=1,JR), (W(I),I=1,JR),H,(A1(I),I=1,N),
         (B1(J), J=1,N)
      OUTPUT 42,22,N,M,JR,IH,L
      OUTPUT 42,23,(XL(J),J=1,JR),(W(I),I=1,JR),C,H,(A1(I),I=1,N),
         (B1(J), J=1, N)
      INPUT 41,1001, IH1, L1, L2
      INPUT 41,1002,H1,S,S1
      INPUT 41,1000, HOLAST, HOLBL, HOLDOT
      RETURN
 1000 FORMAT(3A1)
 1CO1 FORMAT(110)
 1CO2 FORMAT(F10.3)
   20 FORMAT(5110)
   21 FORMAT(E20.8)
   22 FORMAT(1H15110)
   23 FORMAT(1H E20.8)
      ENC
       LIST
       LABEL
C649503
      COMPUTE RETURN G AS FUNCTION OF Y
      SUBROUTINE RETG
      COMMON N.M.JR.H.IH.L.Y.XL.W.A.P.G.NI.A1.B1
     X , IH1, L1, L2, H1, S, S1, HOLAST, HOLBL, HOLDOT
      DIMENSION XL(20), W(20), A(20), P(20), F(20), X(20), A1(10), B1(10)
```

```
IN=INX
      G=XNI+SCRTF(Y+1.)
      RETURN
      END
       LIST
       LABEL
C649504
      NORMALIZED POLYNOMIALS P(K) AS FUNCTIONS OF Y
      SUBROUTINE POLY
      COMMON N.M.JR.H.IH,L.Y.XL.W.A.P.G.NI.A1.BL
     X , IH1, L1, L2, H1, S, S1, HOLAST, HOLBL, HOLDOT
      DIMENSICH XL(20), W(20), A(20), P(20), F(20), X(20), A1(10), B1(10)
      P(2) = SQRTF(3.) +(2.+Y-1.)
      M1=M-2
      DO 40 I=1,M1
      BI=I
      D=SQRTF(2.+BI+3.)/(BI+1.)
      E=SQRTF(2.+81+1.)
      B=BI/SQRTF(2.*BI-1.)
   40 P(I+2)=D+(E+(2.+Y-1.)+P(I+1)-B+P(I))
      RETURN
      END
       LIST
       LABEL
C649505
      COMPUTE AND OUTPUT TOTAL RETURN FUNCTION AT STAGE NI
      SUBROUTINE OUT
      COMMON N.M.JR.H.IH.L.Y.XL.W.A.P.G.NI.A1.B1
     X , IH1, L1, L2, H1, S, S1, HOLAST, HOLBL, HOLDOT
      DIMENSION XL(20), H(20), A(20), P(20), F(20), X(20), A1(10), B1(10)
      DIMENSION Z(110)
      OUTPUT 42,75,NI,JR,M
                                                            M=13)
                     ITERATIVE STAGE
                                         15.5H
                                                 R=13,5H
   75 FORMAT(20H1
       [F(L-1)13,18,18
   13 OUTPUT 42,774
                                              RETURN
  774 FORMAT(40HO
                      RESOURCE
      DO 72 19=1.1H
      X1=19
       Y=(X[-1.) +
      CALL POLY
      SUM=0.
      DO 73 K=1.M
   73 SUM=SUM+A(K)+P(K)
      FF=SUM
      DUTPUT 42.74.Y.FF
   74 FORMAT(1H F15.3,E20.8)
   72 CONTINUE
       GO TO 71
   18 CONTINUE
        CC 772 19=1, IH1
       X1=19
       Y=(X[-1.)+H
       CALL POLY
       SUM=0.
       DO 773 K=1.F
   773 SUM=SUM+A(K)+P(K)
       FF=SUM
       N5=FF+S+S1
       00 1003 15=1,110
```

)

1003 Z(I5)=HOLBL DO 1004 L5=1,L2 IF(I9-L1*L5-1)76,77,76 77 00 78 16=1,110 Z(I6)=HOLDOT 78 CONTINUE 76 CONTINUE 1004 CONTINUE DC 1005 I5=1,110,10 1005 Z(I5)=HOLDOT Z(N5)=HCLAST OUTPUT 42,1006, Y, (2(17), 17=1,105) 772 CONTINUE 71 CONTINUE RETURN 1006 FORMAT(1H F10.3,109A1)

FND

```
Appendix B
2 DIMENSIONAL ALLOCATION VIA DYNAMIC PROGRAMMING. APPROXIMATIONS
C
       LIST
       LABEL
C649500
      COMMON NoMoJRONIOXOYOGOFIOWOVOXLOYLOAOPOPHOFOBOCOIHOHOAIOBIOCIODI
      DIMENSION W(20) + V(20) + XL(20) + YL(20) + A(20+20) + P(20) + PH(20) + F(20+20)
           ,B(3),C(3),SUM(20),SUM2(20),PX(20,20),PY(20,20)
     X
        .B1(20).A1(20).C1(20).D1(20)
    1 CALL INPUT
      IF(N-100)2.3.3
C
      STAGE 1
    2 NI=1
      OUTPUT 42.60.NI.JR.M
   60 FORMAT(20H1
                     ITERATIVE STAGE
                                        15.5H
                                                 R=13.5H
                                                            M=13)
      OUTPUT 42,54
   54 FORMAT(90HO
                                                  Y
                          YDEC
     X
                XDEC
      DO 4 I=1.JR
      X=XL(I)
      CALL POLY
      DO 600 K=1.M
  600 PX(K,I)=P(K)
      DO 41 J=1.JR
      Y=YL(J)
      CALL PHOLY
      DO 601 L=1.M
  601 PY(L,J)=PH(L)
      CALL RETG
      F(1,J)=G
      XDEC=XL(I)
      YDEC=YL(J)
      OUTPUT 42,61,X,Y,F(I,J),XDEC,YDEC
   61 FORMAT(1H 3E20.6.2F10.3)
   41 CONTINUE
    4 CONTINUE
      DO 5 K=1.M
      DO 51 L=1.M
      DO 52 I=1.JR
      SUM(1)=0.
      DO 402 J=1.JR
 402 SUM(I)=SUM(I)+V(J)*F(I,J)*PY(L,J)
   52 CONTINUE
      SUM1 = 0 .
      DO 70 I=1.JR
   70 SUM1=SUM1+W(I) *PX(K.I) *SUM(I)
      A(K.L)=SUM1
  51 CONTINUE
   5 CONTINUE
      OUTPUT 42,50
                      COEFFICIENTS A(K.L)
  50 FORMAT(30H1
                                                        )
      OUTPUT 42,53,((A(I,J),I=1,M),J=1,M)
   53 FORMAT(1H 20.6)
      OUTPUT 42.556
      DO 55 I=1.3
      X=B(1)
```

```
Y=C(1)
    CALL OUT
 55 CONTINUE
    STAGE NI GREATER THAN 1
    DO 6 N2=2.N
    NI=NI+1
    OUTPUT 42,60,NI,JR,M
    OUTPUT 42.54
    DO 8 I=1,JR
    DO 80 J=1.JR
    x=0.
    Y=0.
    CALL RETG
    X=XL(1)-0.
    Y=YL(J)-0.
    CALL TOTRET
    F(I,J)=F1
    XDEC=0.
    YDEC=0.
    DO 9 Il=1.IH
    XI = I1
    IF((XI-1.)*H-A1(I))9,99,99
99 CONTINUE
    IF((XI-1.)*H-B1(I))991,991,11
991 CONTINUE
    IF((XI-1.)*H-XL(I))10,10,11
 10 CONTINUE
    DO 91 J1=1,IH
    XJ=J1
    IF((XJ-1.)*H-C1(J))91,992,992
992 CONTINUE
    IF((XJ-1.)*H-D1(J))993,993,9
993 CONTINUE
    IF((XJ-1.)*H-YL(J))100,100,9
100 X=(XI-1.)*H
    Y=(XJ-1.)*H
    CALL RETG
    X=XL(I)-(XI-1.)+H
    Y=YL(J)-(XJ-1.)*H
    CALL TOTRET
    F2=F1
    IF(F2-F(I,J))9:,91,15
15 F(1,J)=F2
    XDEC=(X1-1.)*H
    YDEC=(XJ-1.)+H
91 CONTINUE
 9 CONTINUE
11 CONTINUE
   OUTPUT 42,61, XL(I),YL(J),F(I,J),XDEC,YDEC
80 CONTINUE
 8 CONTINUE
   DO 400 K=1.M
   DO 401 L=1.M
   DO 502 I=1.JR
```

```
SUM(1)=0.
      DO 500 J=1,JR
  500 SUM(I)=SUM(I)+V(J)*F(I,J)*PY(L,J)
  502 CONTINUE
      SUM1=0.
      DO 501 I=1,JR
  501 SUM1=SUM1+W(I) *SUM(I) *PX(K+I)
      A(K.L)=SUM1
  401 CONTINUE
  400 CONTINUE
      OUTPUT 42.50
      OUTPUT 42,53,((A(I,J),I=1,M),J=1,M)
      OUTPUT 42.556
  556 FORMAT(54HO TEST
                                                                           )
      DO 555 I=1.3
      X=B(I)
      Y=C(I)
      CALL OUT
  555 CONTINUE
    6 CONTINUE
      GO TO 1
    3 CALL EXIT
      STOP
      END
       LIST
       LABEL
C649501
      INPUT ALL CONSTANTS
      SUBROUTINE INPUT
      COMMON N.M.JR.NI.X.Y.G.FI.W.V.XL.YL.A.P.PH.F.B.C.IH.H.A1.B1.C1.D1
      DIMENSION W(20), V(20), XL(20), YL(20), A(20,20), P(20), PH(20), F(20,20)
          .B(3).C(3).SUM(20).SUM2(20).PX(20.20).PY(20.20)
     X
        ,B1(2C),A1(20),C1(20),D1(20)
      INPUT 41.20.N.M.JR.IH
      INPUT 41,21,(XL(J),J=1,JR),(W(I),I=1,JR),(YL(J),J=1,JR),
     X
                   (V(I), I=1,JR), H, (B(I), I=1,3), (C(J), J=1,3)
      OUTPUT42.22.N.A.JR.IH
      OUTPUT42,23,(XL(J),J=1,JR),(W(I),I=1,JR),(B(I),I=1,3),H
      INPUT 41,233, (A1(I), I=1, JR), (B1(I), I=1, JR), (C1(I), I=1, JR),
         (D1(I) • I=1 • JR)
  233 FORMAT(F10.3)
   20 FORMAT(4110)
   21 FORMAT(E20.8)
   22 FORMAT(1H14I10)
   23 FORMAT(1H E20.8)
      RETURN
      END
       LIST
       LABEL
C649502
      COMPUTE RETURN G AS FUNCTION OF X AND Y
      SUBROUTINE RETG
      COMMON NoMOJRONIOXOYOGOFIOWOVOXLOYLOAOPOPHOFOBOCOIHOHOAIOBIOCIODI
      DIMENSION W(20), V(20), XL(20), YL(20), A(20,20), P(20), PH(20), F(20,20)
```

```
.B(3).C(3).SUM(20).SUM2(20).PX(20,20).PY(20,20)
      X
        •B1(20)•A1(20)•C1(20)•D1(20)
       XNI=NI
       G=SQRTF(2.*XNI-1.)*SQRTF(SQRTF((X+.1)*(Y+.1)))
       RETURN
       END
        LIST
        LABEL
C649503
       NORMALIZED POLYNOMIALS P(K)AS FUNCTIONS OF X
       SUBROUTINE POLY
       COMMON N.M.JR.NI.X.Y.G.FI.W.V.XL.YL.A.P.PH.F.B.C.IH.H.AI.BI.CI.DI
       DIMENSION W(20) + V(20) + XL(20) + YL(20) + A(20+20) + P(20) + PH(20) + F(20+20)
           .B(3).C(3).SUM(20).SUM2(20).PX(20,20).PY(20,20)
         •B1(20)•A1(20)•C1(20)•D1(20)
      P(1)=1.
      P(2)=SQRTF(3.)*(2.*X-1.)
       M1 = M - 2
      DO 40 I1=1.M1
       BI = I1
      D=SQRTF(2.*BI+3.)/(BI+1.)
      E=SQRTF(2.#81+1.)
       T=BI/SQRTF(2.*EI-1.)
   40 P(I1+2)=D*(E*(2.*X-1.)*P(I1+1)-T*P(I1))
       RETURN
      END
       LIST
       LABEL
C649504
      NORMALIZED POLYNOMIALS PH(K) AS FUNCTIONS OF Y
      SUBROUTINE PHOLY
      COMMON NoMoJRoNIoXoYoGoFlowoVoXLoYLoAoPoPhoFoBoCoIHoHoAloBloCloDl
      DIMENSION W(20), V(20), XL(20), YL(20), A(20,20), P(20), PH(20), F(20,20)
     X
           .B(3),C(3),SUM(20),SUM2(20),PX(20,20),PY(20,20)
        •B1(20)•A1(20)•C1(20)•D1(20)
      PH(1)=1.
      PH(2)=SQRTF(3.)*(2.*Y-1.)
      M1=M-2
      DO 90 I1=1.M1
      31=11
      D=SGRTF(2.*BI+3.)/(BI+1.)
      E=SQRTF(2.*BI+1.)
      T=BI/SQRTF(2.*BI-1.)
   90 PH(I1+2)=D*(E*(2.*Y-1.)*PH(I1+1)-T*PH(I1))
      RETURN
      END
       LIST
       LABEL
C649505
      COMPUTE TOTAL RETURN AS FUNCTION OF X.Y
      SUBROUTINE TOTRET
      COMMON NoMoJRoNIoxoyoGoFlowovoxLoyLoAoPoPhoFoBoCoIHoHoAloBloCloDl
      DIMENSION W(20) + V(20) + XL(20) + YL(20) + A(20+20) + P(20) + PH(20) + F(20+20)
          .B(3).C(3).SUM(20).SUM2(20).PX(20.20).PY(20.20)
```

```
x .B1(20),A1(20),C1(20),D1(20)
      CALL POLY
      CALL PHOLY
      DO 81 IL=1.M
      SUM2(IL)=0.
      DO 83 IK=1.M
   83 SUM2(IL)=SUM2(IL)+P(IK)*A(IK,IL)
   81 CONTINUE
      SUM3=0.
      DO 84 IL=1.M
   84 SUM3=SUM3+PH(IL)*SUM2(IL)
      F1=G+SUM3
      RETURN
      END
       LIST
       LABEL
C649506
      TOTAL RETURN FROM POLYNOMIAL APPROXIMATION FUNCTION OF X+Y
      SUBROUTINE OUT
      COMMON N.M.JR.NI.X.Y.G.F1.W.V.XL.YL.A.P.PH.F.B.C.IH.H.A1.B1.C1.D1
      DIMENSION W(20) . V(20) . XL(20) . YL(20) . A(20,20) . P(20) . PH(20) . F(20,20)
          ,B(3),C(3),SUM(20),SUM2(20),PX(20,20),PY(20,20)
        .B1(20).A1(20).C1(20).D1(20)
      CALL POLY
      CALL PHOLY
      DO 281 IL=1.M
      SUM2(IL)=0.
      DO 283 IK=1.M
  283 SUM2(IL) = SUM2(IL) +P(IK) *A(IK+IL)
  281 CONTINUE
      SUM3=0.
      DO 284 IL=1.M
  284 SUM3=SUM3+PH(IL) *SUM2(IL)
      FF=SUM3
      ΟUTPUT 42.261.λ.Υ.FF
      RETURN
  261 FORMAT(1H 3E20.6)
      END
                                                                       000255
```

••••

000255

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